### Compulsory 2

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2022-05-06

```
Task 1
```

```
Problem 1.1
library(bootstrap)
library(boot)
library(ggplot2)
library(data.table)
library(Rcpp)
library(VGAM)
## Loading required package: stats4
## Loading required package: splines
##
## Attaching package: 'VGAM'
## The following objects are masked from 'package:boot':
##
       logit, simplex
##
## The following object is masked from 'package:bootstrap':
##
##
       hormone
library(MASS)
# Presenting the data
data <- scor
data.open <- c("mec", "vec")</pre>
data.closed <- c("alg", "ana", "sta")</pre>
all.sets <- c(data.open, data.closed)</pre>
#constructing groups that are to be correlated as functon for the boot package
correlating.groups <- function(data, ind){</pre>
                     mek <- data[ind, 1]</pre>
                     vec <- data[ind, 2]</pre>
                     alg <- data[ind, 3]</pre>
                     ana <- data[ind, 4]
                     sta <- data[ind, 5]</pre>
        c(cor(mek,vec), cor(alg,ana), cor(alg, sta), cor(ana,sta))
```

```
}
boot(data = data, statistic = correlating.groups, R = 200)
## ORDINARY NONPARAMETRIC BOOTSTRAP
##
##
## Call:
## boot(data = data, statistic = correlating.groups, R = 200)
##
##
## Bootstrap Statistics :
                       bias std. error
       original
## t1* 0.5534052 0.0015970223 0.07349188
## t2* 0.7108059 0.0032297071 0.05060925
## t3* 0.6647357 -0.0008492849 0.05792702
## t4* 0.6071743 0.0096818818 0.06458413
```

### problem 1.2

```
theta.hat <- function(data, ind){</pre>
                     mek <- data[ind, 1]</pre>
                     vec <- data[ind, 2]</pre>
                     alg <- data[ind, 3]</pre>
                     ana <- data[ind, 4]</pre>
                     sta <- data[ind, 5]</pre>
                     data <- cbind(mek, vec, alg, ana, sta)</pre>
                     lambda <- eigen(cov(data))$values</pre>
                     lambda[1]/sum(lambda)
theta.boot <- boot(data = data, statistic = theta.hat, R=200, mle = eigen(cov(data)))
theta.boot
## ORDINARY NONPARAMETRIC BOOTSTRAP
##
##
## Call:
## boot(data = data, statistic = theta.hat, R = 200, mle = eigen(cov(data)))
##
## Bootstrap Statistics :
##
       original
                      bias std. error
## t1* 0.619115 0.004075202 0.0474656
```

### problem 1.3

```
print(boot.ci(theta.boot,type = "perc"))

## BOOTSTRAP CONFIDENCE INTERVAL CALCULATIONS
## Based on 200 bootstrap replicates
##

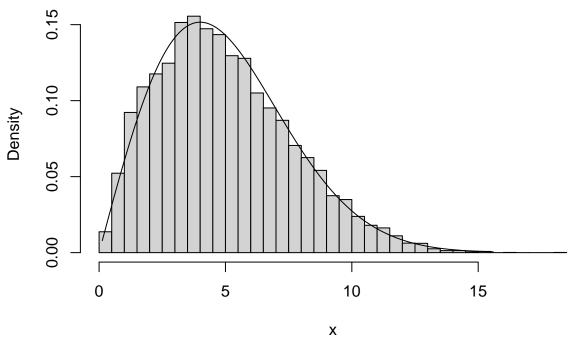
## CALL:
## boot.ci(boot.out = theta.boot, type = "perc")
##

## Intervals:
## Level Percentile
## 95% ( 0.5181,  0.7093 )
## Calculations and Intervals on Original Scale
## Some percentile intervals may be unstable
```

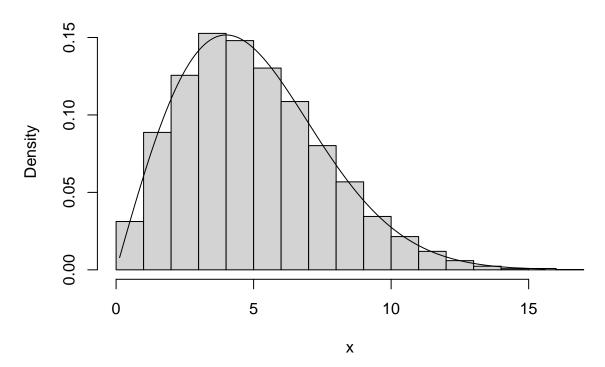
 $\underline{\text{Task 2}}$  problem 2.1 MH-Sampler with rayleigh distribution as proposal distribution

```
set.seed(123)
#Proposal distribution (Rayleigh)
f <- function(x, sigma) {</pre>
 if (any(x < 0)) return (0)
 stopifnot(sigma > 0)
 return((x / sigma^2) * exp(-x^2 / (2*sigma^2)))
}
Nsim <- 20000 #total simulations
burnin = 5000 #sample to burn
sigma <- 4
x <- numeric(Nsim)</pre>
x[1] <- rgamma(1, rate = 1, shape = 1) # start value to simplify index in loop
k <- 0 ## conter for rejects ##
u <- runif(Nsim)</pre>
for (i in 2:Nsim) {
  xt \leftarrow x[i-1]
 y <- rgamma(1, shape = xt, rate = 1)
 num <- f(y, sigma) * dgamma(xt, rate = 1, shape = y) #the numerator of r(xt, y)
  den <- f(xt, sigma) * dgamma(y, rate = 1, shape = xt) #the denominator of <math>r(xt, y)
  if (u[i] \le num/den) x[i] \le y else {
   x[i] <- xt
    k <- k+1 # Number of y being rejected
  }
}
index<-seq(10,Nsim,by=10) ## take index every 10 steps</pre>
#sampled distribution from MH (and library for compartion)
sampled_ray_dist <- x[burnin:Nsim]</pre>
raylib <- rrayleigh(Nsim+1-burnin, scale = sigma)</pre>
#quantile function
a <- ppoints(1000)
QR \leftarrow sigma * sqrt(-2 * log(1 - a))
#Hists
hist(sampled_ray_dist, breaks="scott",
     main="MH-sampler and QF of Raylight", xlab="x", freq=FALSE)
lines(QR, f(QR, 4))
```

# MH-sampler and QF of Raylight



Rayl. dist. from lib. sampler and QF.



### Problem 2.2

Vi Benytter normal fordelingen som symetrisk distribusjon, med  $\sigma = (0.05, 0.5, 2, 16)$ , og på gøy lager vi en sidekick med laplace

```
rw.Metropolis.normal <- function(n, sigma, x0, N) {</pre>
                                     x <- numeric(N)</pre>
                                    x[1] < -x0
                                    u <- runif(N)
                                    k <- 0
                                              for (i in 2:N) {
                                                                y \leftarrow rnorm(1, x[i-1], sigma) \#rlaplace(n=1, location=x[i-1], scale = si
                                                                                  if (u[i] <= (dt(y, n) / dt(x[i-1], n)))</pre>
                                                                                  x[i] \leftarrow y else {
                                                                                  x[i] \leftarrow x[i-1]
                                                                                                    k < - k + 1
                                                                                                     }
                                                                }
                                    return(list(x=x, k=k))
}
rw.Metropolis.laplace <- function(n, sigma, x0, N) {</pre>
                                    x <- numeric(N)
                                    x[1] < -x0
                                    u <- runif(N)
                                    k <- 0
                                             for (i in 2:N) {
                                                                y <- rlaplace(n=1, location=x[i-1], scale = sigma) #
                                                                                   if (u[i] <= (dt(y, n) / dt(x[i-1], n)))</pre>
                                                                                  x[i] <- y else {
                                                                                  x[i] \leftarrow x[i-1]
                                                                                                     k < - k + 1
                                                                                                     }
                                    return(list(x=x, k=k))
```

Random Walk for samplet fra normal fordelingen:

```
n <- 4 #degrees of freedom for target Student t dist.
N <- 2000
sigma <- c(.05, .5, 2, 16)
x0 <- 25
rw1.n <- rw.Metropolis.normal(n, sigma[1], x0, N)
rw2.n <- rw.Metropolis.normal(n, sigma[2], x0, N)
rw3.n <- rw.Metropolis.normal(n, sigma[3], x0, N)
rw4.n <- rw.Metropolis.normal(n, sigma[4], x0, N)
# Result from the random walks given sigma vector
print(c(rw1.n$k, rw2.n$k, rw3.n$k, rw4.n$k))</pre>
```

### **##** [1] 7 225 911 1804

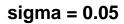
#### ## [1] 0.000 0.099 0.116 0.167

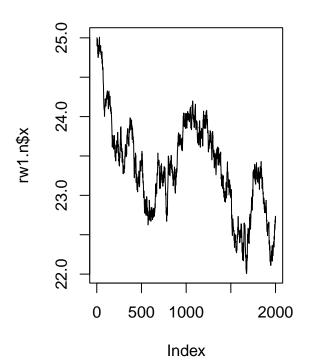
Vi ser at både  $\sigma=2$  og  $\sigma=16$  har en lik andelen innenfor intervallet. For å få en illustrasjon på hvilke  $\sigma$  som konvergerer plotter vi resulktatet:

```
#Plot of sigma1 and sigma2
par(mfrow=c(1,2))

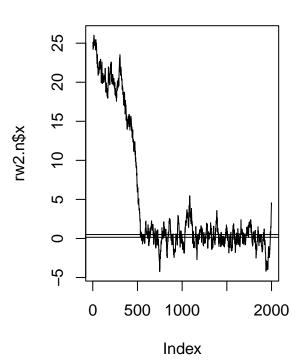
plot(rw1.n$x, type = "l", main="sigma = 0.05")
abline(h=c(0.15,0.5))

plot(rw2.n$x, type = "l", main="sigma = 0.5")
abline(h=c(0.15,0.5))
```



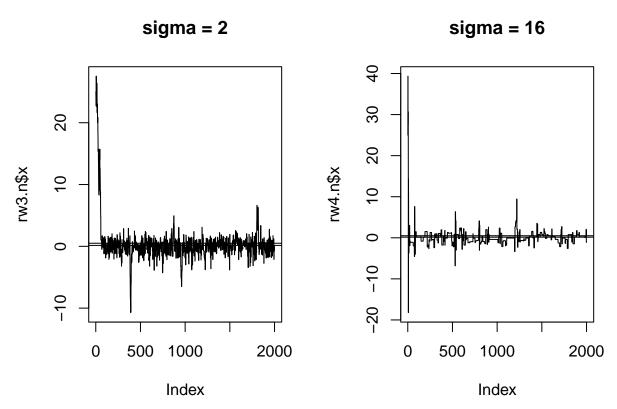


# sigma = 0.5



```
#Plot of sigma1 and sigma2
par(mfrow=c(1,2))
plot(rw3.n$x, type = "l",main="sigma = 2")
abline(h=c(0.15,0.5))

plot(rw4.n$x, type = "l",main="sigma = 16")
abline(h=c(0.15,0.5))
```



En ser på plottet at både  $\sigma = 0.5$ ,  $\sigma = 2$  og  $\sigma = 16$  konvergerer innenfor intervallet, men  $\sigma = 0.5$  trenger en lengre periode med burning før den setler seg i en omegn av intervallet.  $\sigma = 16$  ser noe mindre effektiv ut. Foretrukken  $\sigma$  å optimalisere rundt ville vert  $\sigma = 2$ .

og så til vår lille sidekick:

```
n <- 4 #degrees of freedom for target Student t dist.
N <- 2000
sigma <- c(.05, .5, 2, 16)
x0 <- 25
rw1.1 <- rw.Metropolis.laplace(n, sigma[1], x0, N)
rw2.1 <- rw.Metropolis.laplace(n, sigma[2], x0, N)
rw3.1 <- rw.Metropolis.laplace(n, sigma[3], x0, N)
rw4.1 <- rw.Metropolis.laplace(n, sigma[4], x0, N)
# Result from the random walks given sigma vector
print(c(rw1.1$k, rw2.1$k, rw3.1$k, rw4.1$k))
## [1] 9 327 953 1751</pre>
```

```
#Proportion of rejections within [0.15, 0.5]

print(c((2000 - length(which(rw1.1$x > 0.5)) - length(which(rw1.1$x < 0.15)))/N,

(2000 - length(which(rw2.1$x > 0.5)) - length(which(rw2.1$x < 0.15)))/N,

(2000 - length(which(rw3.1$x > 0.5)) - length(which(rw3.1$x < 0.15)))/N,

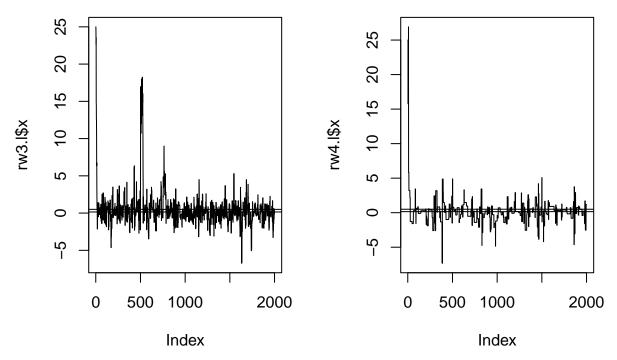
(2000 - length(which(rw4.1$x > 0.5)) - length(which(rw4.1$x < 0.15)))/N)
```

```
## [1] 0.000 0.108 0.119 0.216
```

For å se for hvilke  $\sigma$  som konvergerer

```
#Plot of sigma1 and sigma2
par(mfrow=c(1,2))
```

```
plot(rw1.1$x, type = "1")
abline(h=c(0.15,0.5))
plot(rw2.1$x, type = "1")
abline(h=c(0.15,0.5))
     27.5
                                                    25
                                                    20
     26.5
                                                    15
rw1.I$x
                                                    10
     25.5
                                                    2
                                                    0
     24.5
                                                    -5
           0
                500
                      1000
                                    2000
                                                               500
                                                                    1000
                                                                                  2000
                                                         0
                      Index
                                                                     Index
#Plot of sigma1 and sigma2
par(mfrow=c(1,2))
plot(rw3.1$x, type = "1")
abline(h=c(0.15,0.5))
plot(rw4.1$x, type = "l")
abline(h=c(0.15,0.5))
```

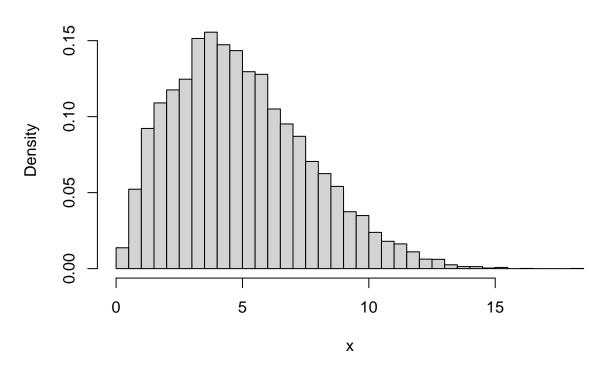


Det må sies å være til forveksling likt i konvergens i tillegg konvergerer  $\sigma=0.05$  noe fortere.

### Problem 2.3

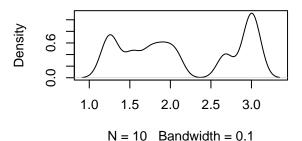
```
set.seed(123)
\#initialize constants and parameters
N \leftarrow 5000 \# length \ of \ chain
burn <- 1000 #burn-in length
X <- matrix(0, N, 2) #the chain, a bivariate sample
rho <- -.9 #correlation
mu_x <- 0
mu_y <- 0
sigma_x <- 1
sigma_y <- 1
s1 <- sqrt(1-rho^2)*sigma_x</pre>
s2 <- sqrt(1-rho^2)*sigma_y</pre>
##### generate the chain #####
X[1, ] <- c(mu_x, mu_y) #initialize</pre>
for (i in 2:N) {
x2 < X[i-1, 2]
m_x \leftarrow mu_x + rho * (y - mu_y) * sigma_x/sigma_y
X[i, 1] <- rnorm(1, m_x, s1)</pre>
x_x < X[i, 1]
m_y \leftarrow mu_y + rho * (x_x - mu_x) * sigma_y/sigma_x
X[i, 2] <- rnorm(1, m_y, s2)
b <- burn + 1
x \leftarrow X[b:N,]
hist(sampled_ray_dist, breaks="scott",
 main="Gibbs Sampler for N(0,1)", xlab="x", freq=FALSE)
```

# Gibbs Sampler for N(0,1)

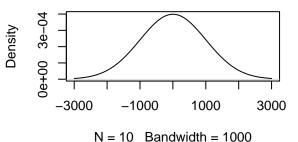


```
Task 3
problem 3.1
p1 <- 0.5
n = 10
data \leftarrow p1 *rnorm(n, 1,1) + (1-p1) * rnorm(n, 3,1)
data1 <- p1 *rgamma(n, rate = 1, shape = 1) + (1-p1) * rgamma(n, rate = 3, shape = 1)
par(mfrow = c(1, 1))
print(density(data))
##
## Call:
## density.default(x = data)
                             Bandwidth 'bw' = 0.4101
## Data: data (10 obs.);
##
##
## Min. :-0.01707
                             :0.001771
                      Min.
## 1st Qu.: 1.05666
                      1st Qu.:0.053027
## Median: 2.13038 Median: 0.283447
## Mean : 2.13038 Mean :0.232491
## 3rd Qu.: 3.20411
                       3rd Qu.:0.374453
          : 4.27783
                               :0.455501
                      Max.
df <- approxfun(density(data)) ##extract the density function##</pre>
xnew \leftarrow seq(0,10,0.2)
n <- length(data)</pre>
h1 \leftarrow 0.9 * min(c(IQR(data)/1.34, sd(data))) * n^(-1/5) #density seems not unimodal#
h2 \leftarrow 1.06 * sd(data) * n^(-1/5)
lets first look at the extream example (high, low):
par(mfrow = c(2, 2))
plot(density(data, bw=.10) , main="extream narrow bw - gaus distr") # "Day of the tentacle"
plot(density(data, bw=1000), main="extream wide bw - gaus distr") #Gleat - "too much
plot(density(data1,bw=.10), main="extream narrow bw - gamma distr") # "Day of the tentacle"
plot(density(data1,bw=1000), main="extream narrow bw - gamma distr") #Gleet - "too much
```

### extream narrow bw - gaus distr

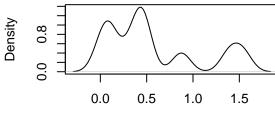


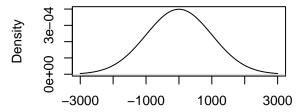
### extream wide bw - gaus distr



### extream narrow bw - gamma distr

### extream narrow bw - gamma distr





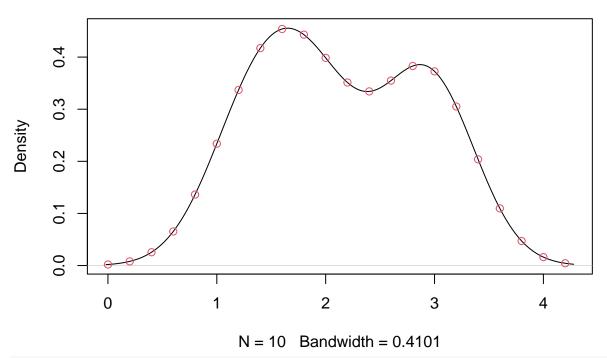
N = 10 Bandwidth = 0.1

N = 10 Bandwidth = 1000

From this we can conclude that choise of bandwith is essential. To narrow you get spikes, to wide you obtain a verry flat, smooth curve converged towards the normal distribution. As we also see the result is the same if we change the kernel distribution. So lets try to fit a better curve:

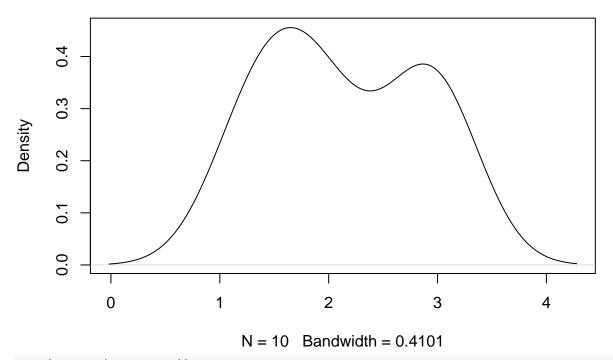
#par(mfrow=c(2,2))plot(density(data)) #The default method applies the Gaussian kernel#3 points(xnew,df(xnew),col=2) #bulky

# density.default(x = data)



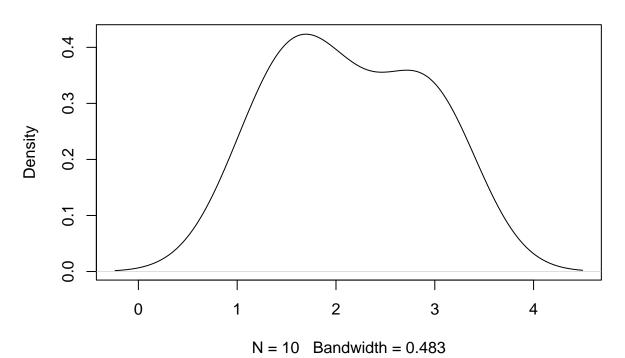
plot(density(data,bw=h1)) #Better, this include the duality of the two distr.

# density.default(x = data, bw = h1)



plot(density(data,bw=h2)) #oversmooth...

# density.default(x = data, bw = h2)



problem 3.2

we have:

$$\mathbf{x} = \{x_1, ..., x_n\} \sim gamma(\theta)$$
  
 $\theta = (r, \lambda)$ 

we first need to solve for the MLE functions for the gamma distributed RVs:

$$f(\mathbf{x}|\theta) = \frac{\lambda^r}{\Gamma(r)} \cdot \mathbf{x}^{r-1} \cdot exp(-\lambda \mathbf{x})$$

$$L(\theta|x_1, ..., x_n) = \prod_{i=1}^n f(\mathbf{x}|\theta)$$

$$= \frac{\lambda^{n \cdot r}}{\Gamma(r)^n} \cdot \sum_i^n x_i^{r-1} \cdot exp(-\lambda \sum_i^n x_i)$$

we obtain the log likeliehood by applying magic

$$l(\theta|x_1,...,x_n) = nr \cdot log(\lambda) - n \cdot log(\Gamma(r)) + (r-1) \cdot \sum_{i=1}^{n} log(x_i) - \lambda \sum_{i=1}^{n} x_i$$

we then take the derivative to find the maximum:

$$\frac{\partial l}{\partial \lambda} = \frac{nr}{\lambda} - \sum_{i=1}^{n} x_i = 0$$

$$\frac{\partial l}{\partial r} = n \cdot \log(\lambda) - \frac{n\Gamma'(r)}{\Gamma(r)} + \sum_{i=1}^{n} \log(x_i) = 0$$

to find  $\hat{\lambda}$  is fairly straight forward:

$$\frac{nr}{\lambda} - \sum_{i=1}^{n} x_i = 0$$

$$\frac{nr}{\lambda} = \sum_{i=1}^{n} x_i$$

$$\hat{\lambda} = \frac{nr}{\sum_{i=1}^{n} x_i} = \frac{r}{n^{-1} \sum_{i=1}^{n} x_i} = \frac{\hat{r}}{\bar{x_n}}$$

to find  $\hat{r}$  is more of a pain since we have the  $\frac{\Gamma'(r)}{\Gamma(r)}$  term also known as digamma:

$$\psi(r) = \frac{\Gamma'(r)}{\Gamma(r)} = \frac{\partial}{\partial r} log(\Gamma(r))$$

to simplify we substitute  $\hat{\lambda}$  into the equation:

$$n \cdot \log(\frac{\hat{r}}{\bar{x_n}}) - \frac{n\Gamma'(\hat{r})}{\Gamma(\hat{r})} + \sum_{i=1}^n \log(x_i) = 0$$
 (1)

$$log(\frac{\hat{r}}{\bar{x_n}}) - \frac{1}{n} \sum_{i=1}^n log(x_i) = \frac{\Gamma'(\hat{r})}{\Gamma(\hat{r})} = \psi(\lambda \bar{x_n})$$
 (2)

for  $\bar{x_n} = \frac{r}{\lambda}$ 

Due to the digamma function the MLE needs to be solved numerically by root solver (in r uniroots), which is presumably a bisection method, since derivative is not needed. Bisection only need a continuous function to converge to one of the existing roots. Bellow the same simulation asfrom "Statistical computing with R by Maria L. Rizzo" page 340 is conducted (with comments on the process)

```
m <- 20000
est <- matrix(0, m, 2)</pre>
n <- 200
r \leftarrow 5
lambda <- 2
# equation (1) above, of which we want to find the root
obj <- function(lambda, xbar, logx.bar) {</pre>
    digamma(lambda * xbar) - logx.bar - log(lambda)
    }
#Do m simulations of a set of x_1, \ldots, x_n
for (i in 1:m) {
    x <- rgamma(n, shape=r, rate=lambda)
    xbar <- mean(x)</pre>
    # using uniroot to solve numericaly
    u <- uniroot(obj, lower = .001, upper = 10e5,
        xbar = xbar, logx.bar = mean(log(x)))
  #columns vector of MLE simulations
  lambda.hat <- u$root</pre>
  r.hat <- xbar * lambda.hat</pre>
  est[i, ] <- c(r.hat, lambda.hat)</pre>
#mean of the MLE simulations
ML <- colMeans(est)</pre>
ML
```

## [1] 5.070865 2.030602